

Sebacic acid, 2,6-dichlorophenyl ethyl ester

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|-----------------------------|---|
| Inchi: | InChI=1S/C18H24Cl2O4/c1-2-23-16(21)12-7-5-3-4-6-8-13-17(22)24-18-14(19)10-9-11-15 |
| InchiKey: | CEQJWUNFXAGIKB-UHFFFAOYSA-N |
| Formula: | C18H24Cl2O4 |
| SMILES: | CCOC(=O)CCCCCCCC(=O)Oc1c(Cl)cccc1Cl |
| Mol. weight [g/mol]: | 375.29 |

Physical Properties

| Property code | Value | Unit | Source |
|---------------|---------|----------------------|----------------|
| gf | -297.87 | kJ/mol | Joback Method |
| hf | -722.34 | kJ/mol | Joback Method |
| hfus | 49.61 | kJ/mol | Joback Method |
| hvap | 86.34 | kJ/mol | Joback Method |
| log10ws | -6.20 | | Crippen Method |
| logp | 5.583 | | Crippen Method |
| mcvol | 280.080 | ml/mol | McGowan Method |
| pc | 1449.04 | kPa | Joback Method |
| rinqol | 2665.00 | | NIST Webbook |
| tb | 875.32 | K | Joback Method |
| tc | 1084.97 | K | Joback Method |
| tf | 548.24 | K | Joback Method |
| vc | 1.081 | m ³ /kmol | Joback Method |

Temperature Dependent Properties

| Property code | Value | Unit | Temperature [K] | Source |
|---------------|-----------|---------|-----------------|---------------|
| cpg | 812.47 | J/molxK | 875.32 | Joback Method |
| cpg | 825.76 | J/molxK | 910.26 | Joback Method |
| cpg | 837.96 | J/molxK | 945.20 | Joback Method |
| cpg | 849.09 | J/molxK | 980.15 | Joback Method |
| cpg | 859.16 | J/molxK | 1015.09 | Joback Method |
| cpg | 868.21 | J/molxK | 1050.03 | Joback Method |
| cpg | 876.23 | J/molxK | 1084.97 | Joback Method |
| dvisc | 0.0004290 | Paxs | 548.24 | Joback Method |
| dvisc | 0.0002609 | Paxs | 602.75 | Joback Method |

| | | | | |
|-------|-----------|------|--------|---------------|
| dvisc | 0.0001723 | Paxs | 657.27 | Joback Method |
| dvisc | 0.0001212 | Paxs | 711.78 | Joback Method |
| dvisc | 0.0000897 | Paxs | 766.29 | Joback Method |
| dvisc | 0.0000691 | Paxs | 820.81 | Joback Method |
| dvisc | 0.0000549 | Paxs | 875.32 | Joback Method |

Sources

| | |
|------------------------|---|
| McGowan Method: | http://link.springer.com/article/10.1007/BF02311772 |
| NIST Webbook: | http://webbook.nist.gov/cgi/cbook.cgi?ID=U354872&Units=SI |
| Crippen Method: | http://pubs.acs.org/doi/abs/10.1021/ci9903071 |
| Crippen Method: | https://www.chemeo.com/doc/models/crippen_log10ws |
| Joback Method: | https://en.wikipedia.org/wiki/Joback_method |

Legend

| | |
|-----------------|---|
| cpg: | Ideal gas heat capacity |
| dvisc: | Dynamic viscosity |
| gf: | Standard Gibbs free energy of formation |
| hf: | Enthalpy of formation at standard conditions |
| hfus: | Enthalpy of fusion at standard conditions |
| hvap: | Enthalpy of vaporization at standard conditions |
| log10ws: | Log10 of Water solubility in mol/l |
| logp: | Octanol/Water partition coefficient |
| mcvol: | McGowan's characteristic volume |
| pc: | Critical Pressure |
| rinpol: | Non-polar retention indices |
| tb: | Normal Boiling Point Temperature |
| tc: | Critical Temperature |
| tf: | Normal melting (fusion) point |
| vc: | Critical Volume |

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